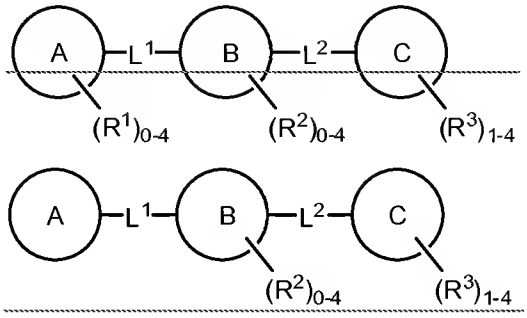
Amendments to the Claims

The following listing of claims will replace all prior versions and listings of claims in the application.

Listing of Claims:

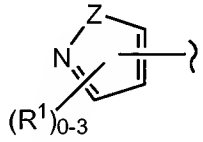
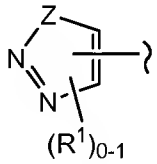
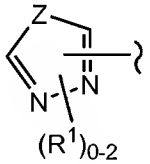
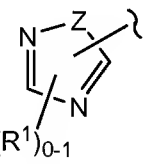
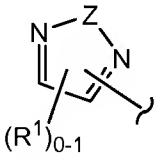
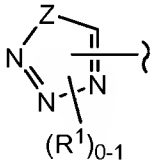
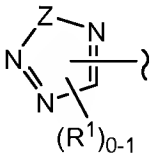
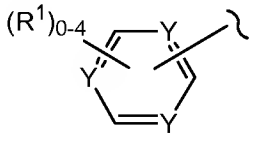
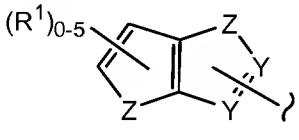
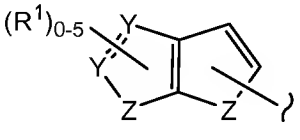
1. (Currently amended) A compound for modulating c-Kit activity according to Formula **I**,

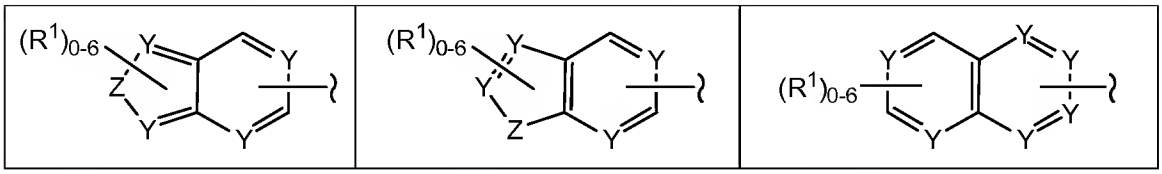


I

or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:



wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and-N(R^7)-, provided that the A ring contains at least one annular N, O, or S;

each R^1 is independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^4, -$S(O)_{0-2}R^4$, -$SO_2N(R^4)R^4$, -$C(=O)N(R^4)R^4$, -$C(=NR^5)N(R^4)R^4$, -$C(=NR^5)R^4$, -$N(R^4)SO_2R^4$, -$N(R^4)C(O)R^4$, -NCO_2R^4, -$C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two adjacent of R^1, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{10};

L^1 is a single bond;

ring B is phenyl;

each R^2 is independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^4, -$N(R^4)R^4$, -$S(O)_{0-2}R^4$, -$SO_2N(R^4)R^4$, -$C(=O)N(R^4)R^4$, -$C(=NR^5)N(R^4)R^4$, -$C(=NR^5)R^4$, -$N(R^4)SO_2R^4$, -$N(R^4)C(O)R^4$, -NCO_2R^4, -$C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two adjacent of R^2, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{15};

L^2 is selected from -N(H)N(H)C(=O)N(H)-, -CH_2N(H)C(=O)N(H)-, -CH_2OC(=O)N(H)-, and -XCH_2C(=O)N(H)-; wherein X is selected from -O-, -$S(O)_{0-2}$-, and -N(R^7)-; and any C-H of L^2 is optionally C-R^{20};

ring C is phenyl;

each R^3 is independently selected from halogen, trihalomethyl, -CN,-NO_2, -OR^4, -$N(R^4)R^4$, -$S(O)_{0-2}R^4$, -$SO_2N(R^4)R^4$, -$C(=O)N(R^4)R^4$, -$C(=NR^5)N(R^4)R^4$, -$C(=NR^5)R^4$, -$N(R^4)SO_2R^4$, -$N(R^4)C(O)R^4$, -NCO_2R^4, -$C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, ~~optionally substituted heterocyclyl,~~ and optionally substituted heterocyclyl C_{1-6}alkyl; provided R^3 is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide,

wherein there exists at least one of R^3 that is halogen or trihalomethyl;

R^4 is selected from -H, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two of R^4, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^5 is selected from -H, -CN, -NO_2, -OR^4, -$S(O)_{0-2}R^4$, -CO_2R^4, optionally substituted C_{1-6}alkyl, optionally substituted C_{1-6}alkenyl, and optionally substituted C_{1-6}alkynyl;

R^7 is selected from -H, optionally substituted C_{1-6}alkyl, -$SO_2N(R^4)R^4$, -CO_2R^4, -$C(=O)N(R^4)R^4$, -$C(=NR^5)N(R^4)R^4$, -$C(=NR^5)R^4$, -$C(=O)R^4$, optionally substituted alkoxy, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl; and

each of R^{10}, each of R^{15}, each of R^{20}, and each of R^{25} is independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^4, -$N(R^4)R^4$, -$S(O)_{0-2}R^4$, -$SO_2N(R^4)R^4$, -CO_2R^4, -$C(=O)N(R^4)R^4$, -$C(=NR^5)N(R^4)R^4$, -$C(=NR^5)R^4$, -$N(R^4)SO_2R^4$, -$N(R^4)C(O)R^4$, -NCO_2R^4, -$C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl

C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

provided:

the compound is not one of:

N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,

N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,

N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide

2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-[3-(trifluoromethyl)phenyl]acetamide

N-[2-chloro-5-(trifluoromethyl) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide or

N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

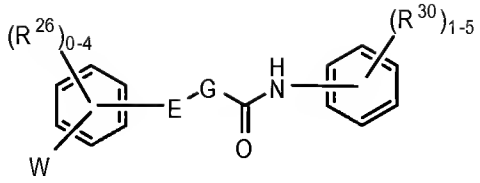
2-12. (Cancelled)

13. (Canceled)

14. (Original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L^2.

15. (Previously presented) The compound according to claim 1, wherein each of R^3 is independently selected from halogen, trihalomethyl, $-OR^4$, $-C(=O)R^4$, and optionally substituted C_{1-6}alkyl.

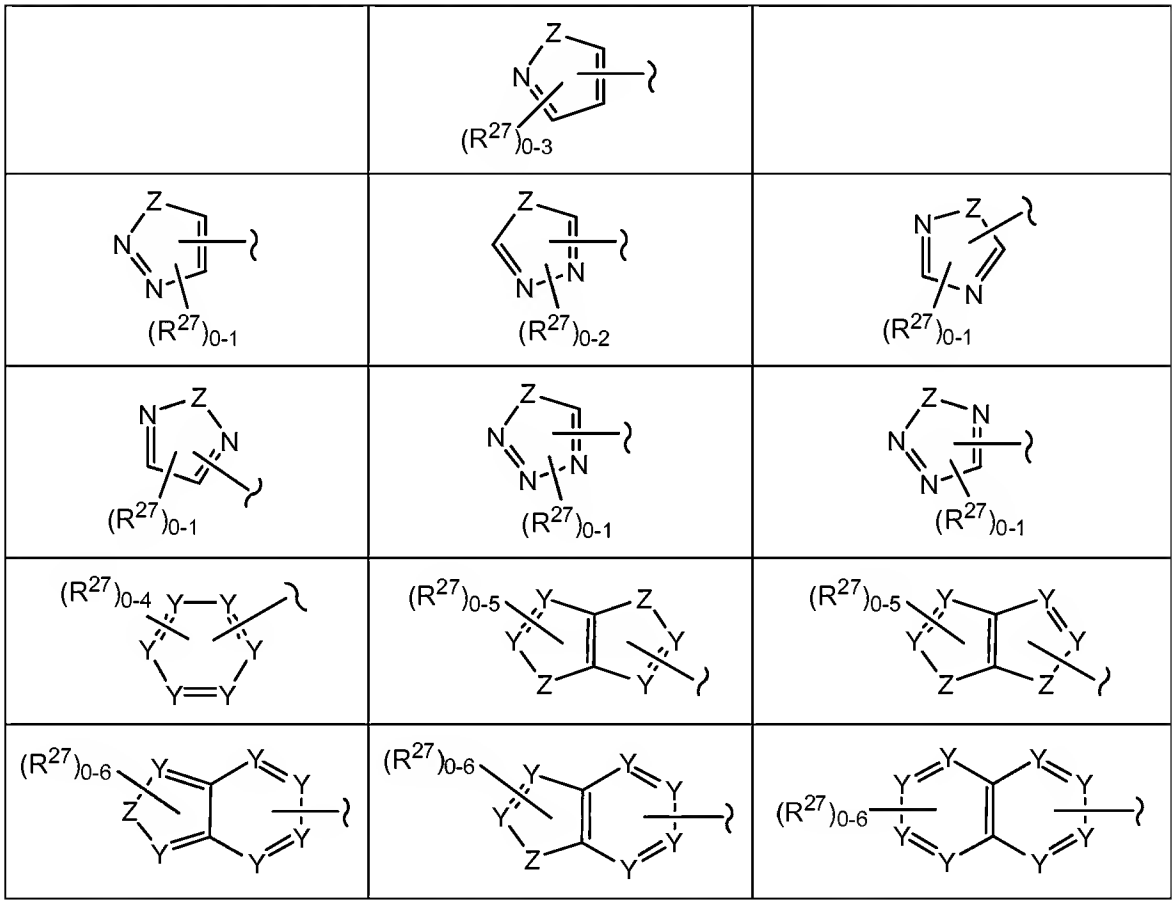
16. (Currently amended) A compound for modulating c-Kit activity according to the following Formula:



II

or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:



each of R^{27} independently selected from halogen, trihalomethyl, -CN, $-NO_2$, $-OR^{55}$, $-S(O)_{0-2}R^{55}$, $-SO_2N(R^{55})R^{55}$, $-C(=O)N(R^{55})R^{55}$, $-C(=NR^{50})N(R^{55})R^{55}$, $-C(=NR^{50})R^{55}$, $-N(R^{55})SO_2R^{55}$, $-N(R^{55})C(O)R^{55}$, $-NCO_2R^{55}$, $-C(=O)R^{55}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -$S(O)_{0-2}$-, and -$N(R^7)$-,

provided that the W ring contains at least one annular annular N, O, or S;

E and G are each independently selected from -O-, -$S(O)_{0-2}$-, -$C(R^{31})R^{32}$-, and -$N(R^{33})$-;

J_1 and J_2 are each independently =C(H)- or =N-;

R^{26} is independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^{40}, -$N(R^{40})R^{40}$, -$S(O)_{0-2}R^{40}$, -$SO_2N(R^{40})R^{40}$, -CO_2R^{40}, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$N(R^{40})SO_2R^{40}$, -$N(R^{40})C(O)R^{40}$, -$NCO_2R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

R^{30} is independently selected from halogen, trihalomethyl, -CN, -NO_2, -OR^{40}, -$N(R^{40})R^{40}$, -$S(O)_{0-2}R^{40}$, -$SO_2N(R^{40})R^{40}$, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$N(R^{40})SO_2R^{40}$, -$N(R^{40})C(O)R^{40}$, -$NCO_2R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, ~~optionally substituted heterocyclyl,~~ and optionally substituted heterocyclyl C_{1-6}alkyl, wherein there exists at least one of R^{30} that is trihalomethyl;

$R^{31}$and R^{32} are each independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^{40}, -$N(R^{40})R^{40}$, -$S(O)_{0-2}R^{40}$, -$SO_2N(R^{40})R^{40}$, -CO_2R^{40}, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$N(R^{40})SO_2R^{40}$, -$N(R^{40})C(O)R^{40}$, -$NCO_2R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

R^{33} is selected from -H, optionally substituted lower alkyl, -$SO_2N(R^{40})R^{40}$, -CO_2R^{40}, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

R^{40} is selected from -H, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two of R^{40}, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^{50} is selected from -H, -CN, $-NO_2$, $-OR^{40}$, $-S(O)_{0-2}R^{40}$, $-CO_2R^{40}$, optionally substituted C_{1-6}alkyl, optionally substituted C_{1-6}alkenyl, and optionally substituted C_{1-6}alkynyl;

R^{55} is selected from -H, optionally substituted C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl; and

two of R^{55}, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

17. (Cancelled)

18. (Currently amended) The compound according to claim 16, wherein R^{30} is selected from halogen, trihalomethyl, $-OR^{40}$, $-N(R^{40})R^{40}$, $-C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, ~~optionally substituted heterocyclyl,~~ and optionally substituted heterocyclyl C_{1-6}alkyl, wherein there exists at least one of R^{30} that is trifluoromethyl.

19. (Cancelled)

20. (Cancelled)

21. (Cancelled)

22. (Cancelled)

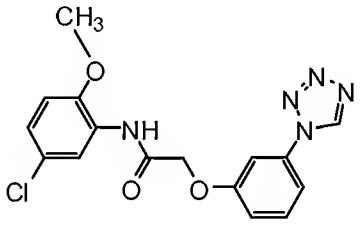
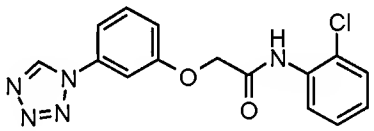
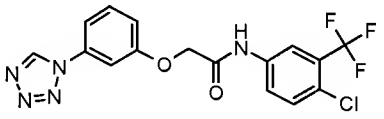
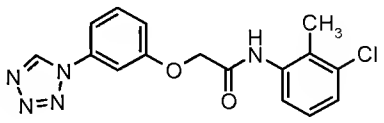
23. (Withdrawn) The compound according to claim 16, wherein E is selected from -O-, $-S(O)_{0-2}-$, and -NH-; and G is $-CH_2-$.

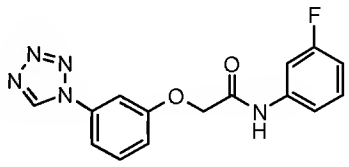
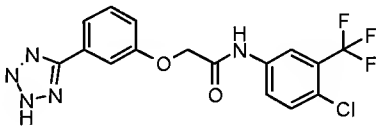
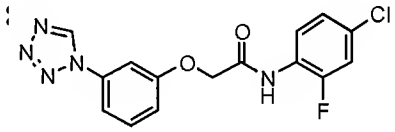
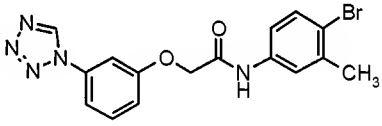
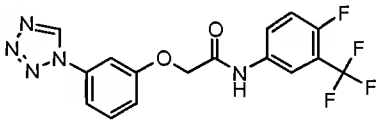
24. (Withdrawn) The compound according to claim 16, wherein E is either $-CH_2-$ or -NH-; and G is selected from -O-, -S-, and -NH-.

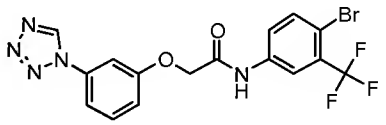
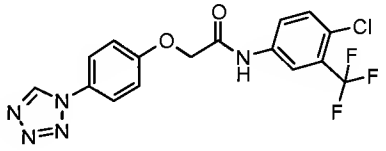
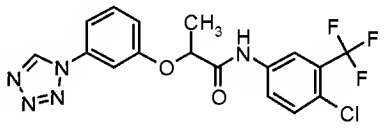
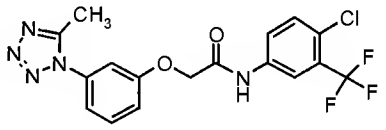
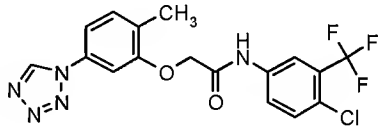
25. (Cancelled)

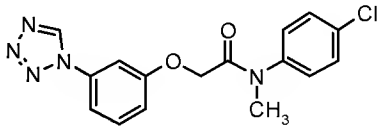
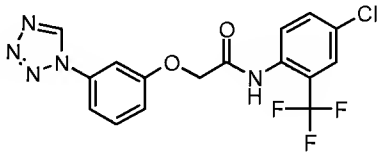
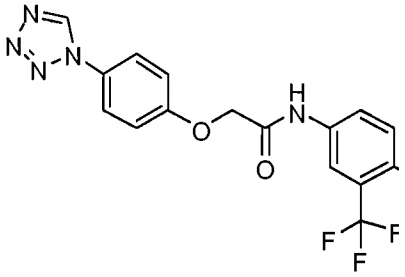
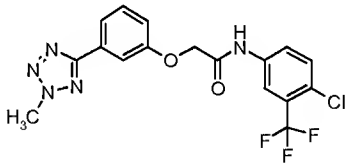
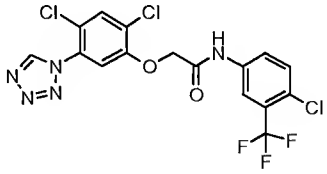
26. (Cancelled)

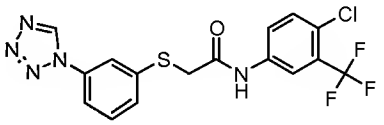
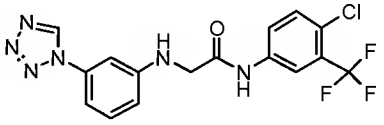
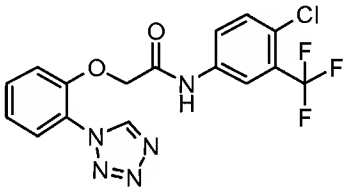
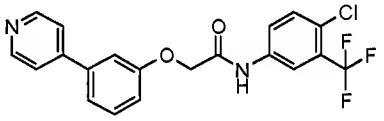
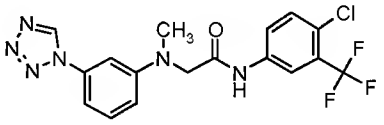
27. (Previously presented) A compound selected from the following Table:

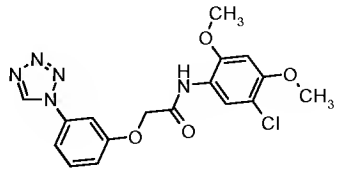
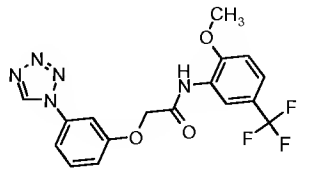
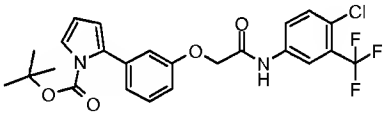
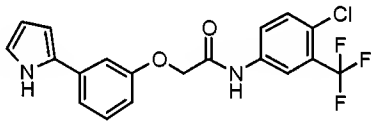
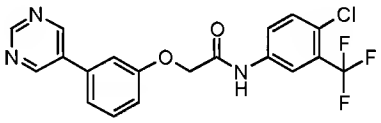
Entry	Name	Structure
1	N-[5-chloro-2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
4	N-(2-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
7	N-(3-chloro-2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

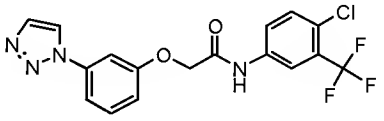
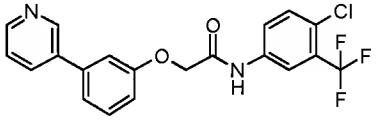
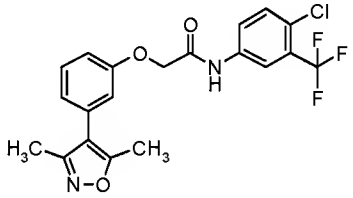
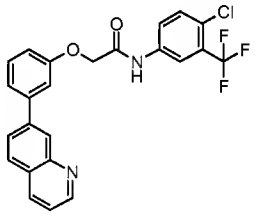
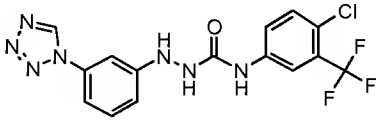
Entry	Name	Structure
8	N-(3-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2H-tetrazol-5-yl)phenyl]oxy}acetamide	
10	N-(4-chloro-2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

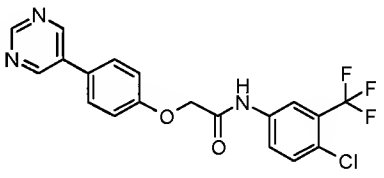
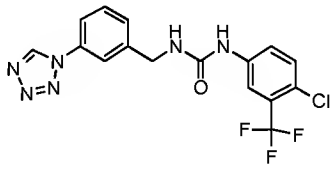
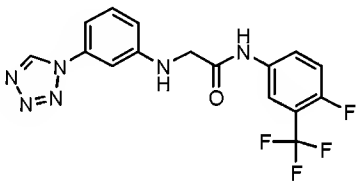
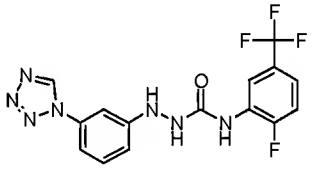
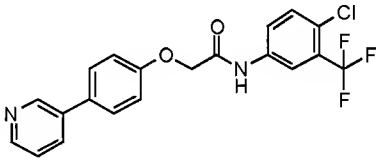
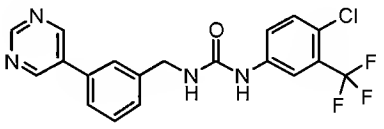
Entry	Name	Structure
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(5-methyl-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-methyl-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

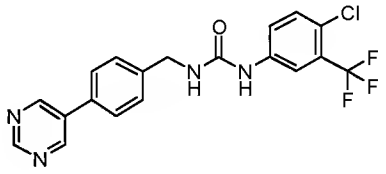
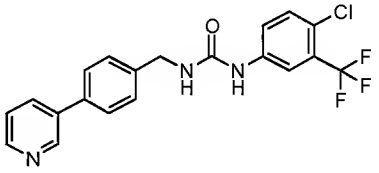
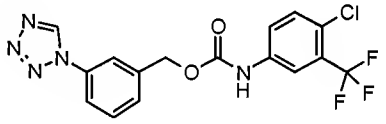
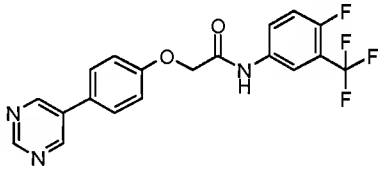
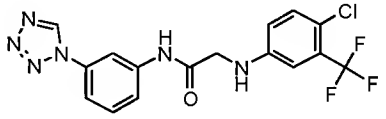
Entry	Name	Structure
19	N-(4-chlorophenyl)-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2-methyl-2H-tetrazol-5-yl)phenyl]oxy}acetamide	
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,4-dichloro-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

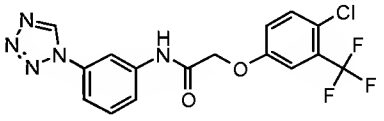
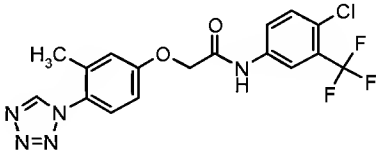
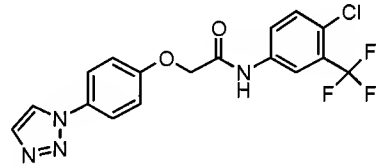
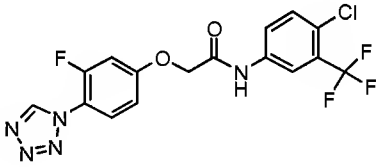
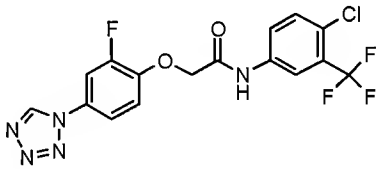
Entry	Name	Structure
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide	
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

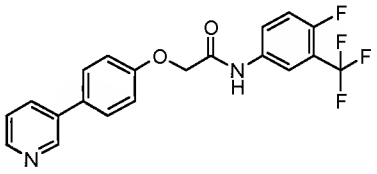
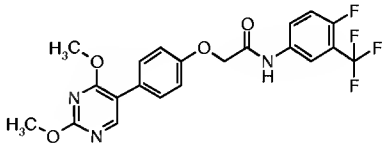
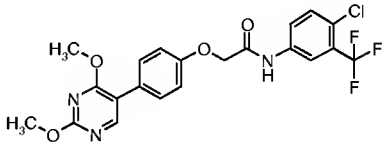
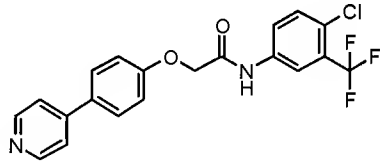
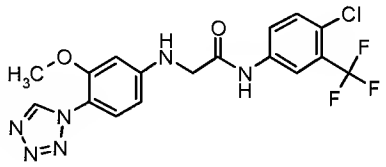
Entry	Name	Structure
49	N-[5-chloro-2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
52	N-[2-(methyloxy)-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
57	1,1-dimethylethyl 2-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate	
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-pyrrol-2-yl)phenyl]oxy}acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyrimidin-5-ylphenyl)oxy]acetamide	

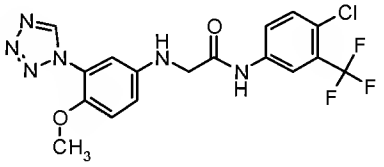
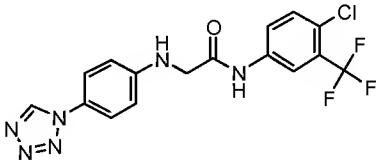
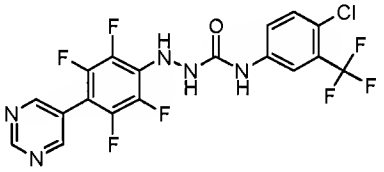
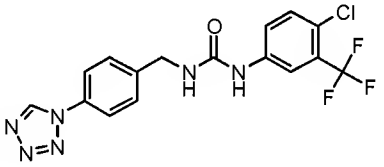
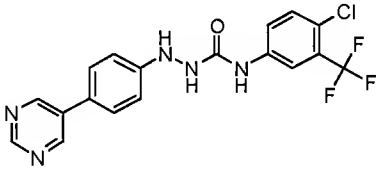
Entry	Name	Structure
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide	
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-3-ylphenyl)oxy]acetamide	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(3,5-dimethylisoxazol-4-yl)phenyl]oxy}acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-quinolin-7-ylphenyl)oxy]acetamide	
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	

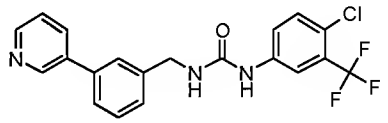
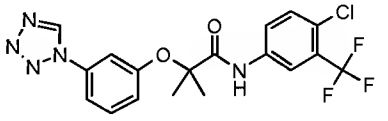
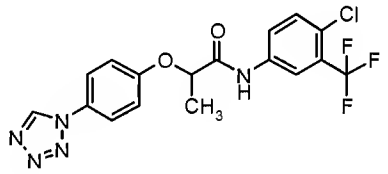
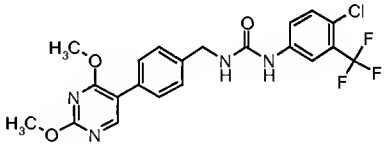
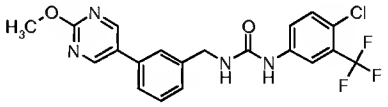
Entry	Name	Structure
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	

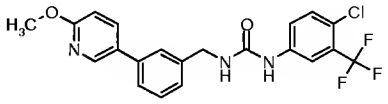
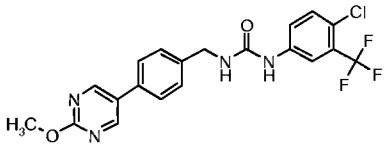
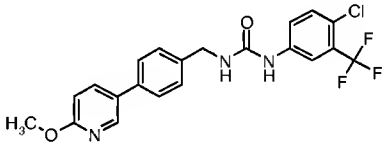
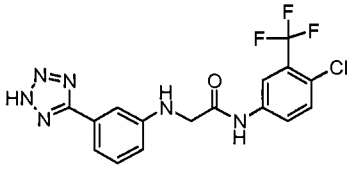
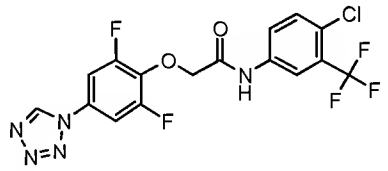
Entry	Name	Structure
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
86	N~2~-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

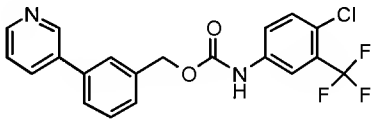
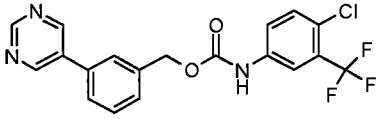
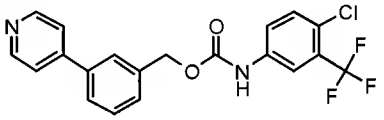
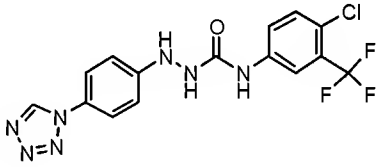
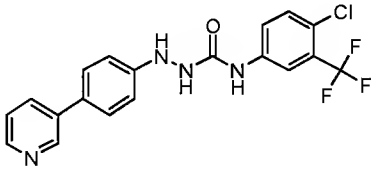
Entry	Name	Structure
87	2-{[4-chloro-3-(trifluoromethyl)phenyl]oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

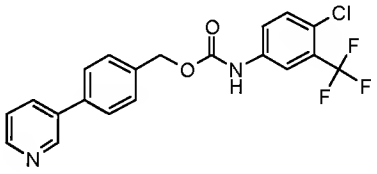
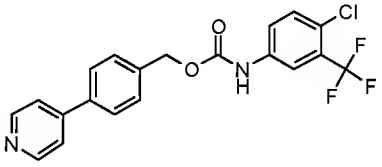
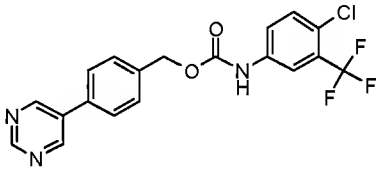
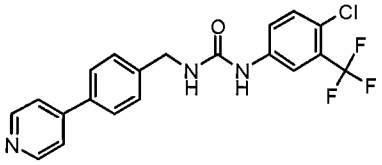
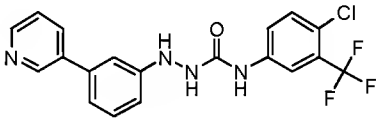
Entry	Name	Structure
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	
96	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	

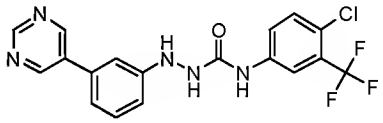
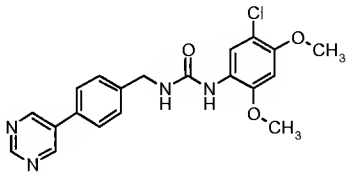
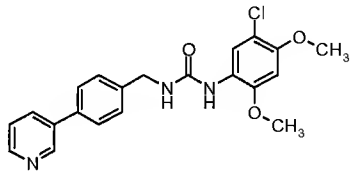
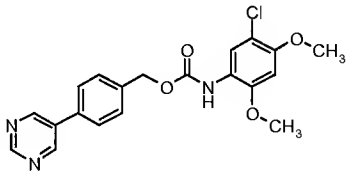
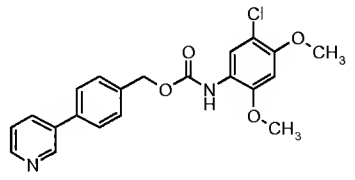
Entry	Name	Structure
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methyloxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-tetrazol-1-yl)phenyl]methyl}urea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	

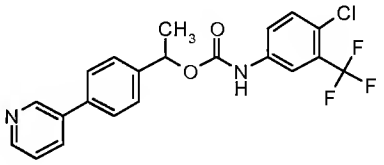
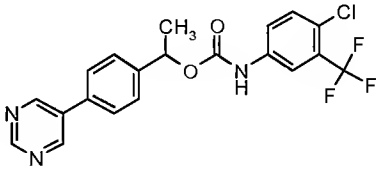
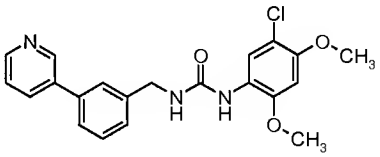
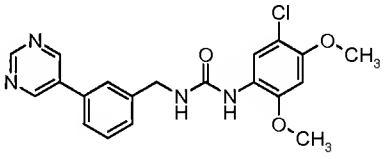
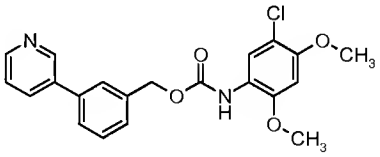
Entry	Name	Structure
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
107	N-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	

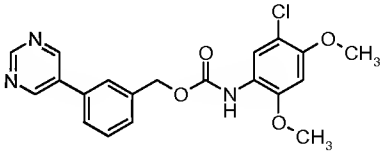
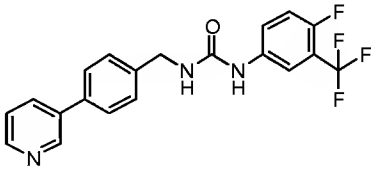
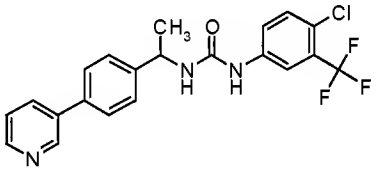
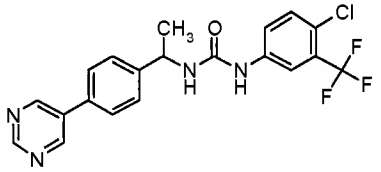
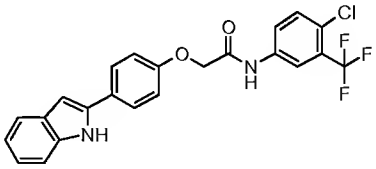
Entry	Name	Structure
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

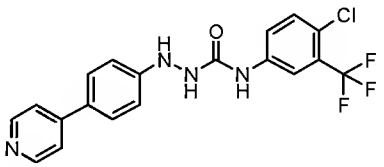
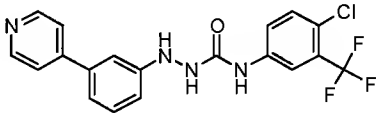
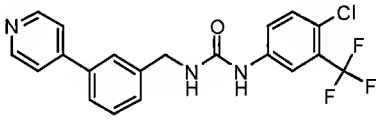
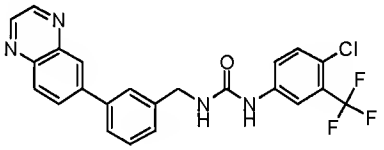
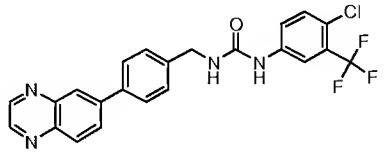
Entry	Name	Structure
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	

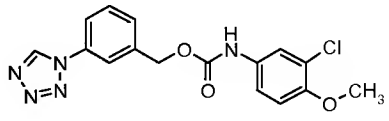
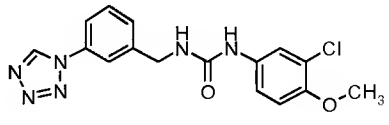
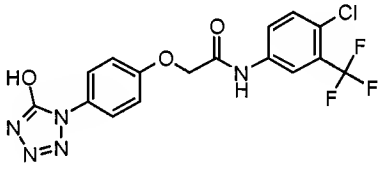
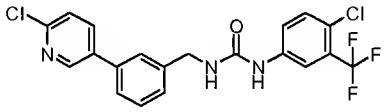
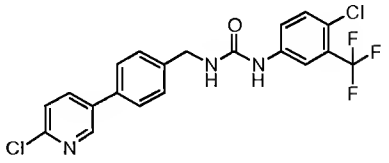
Entry	Name	Structure
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	

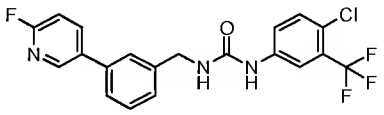
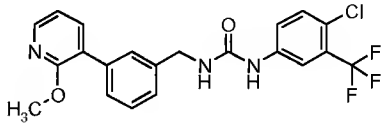
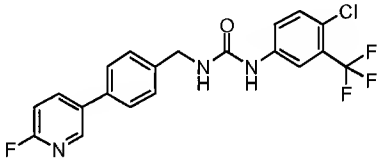
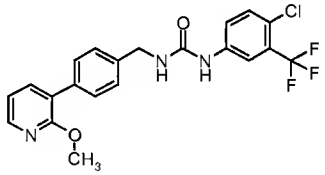
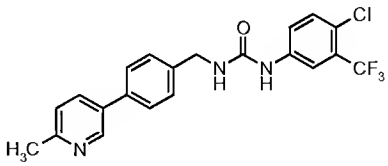
Entry	Name	Structure
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
128	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	

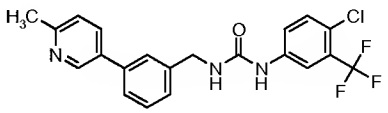
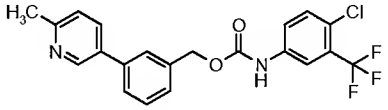
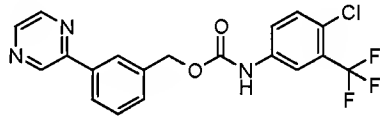
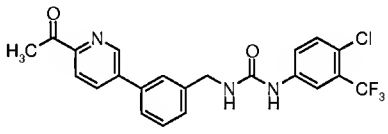
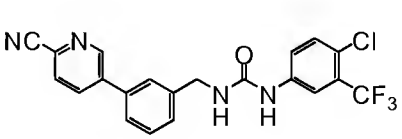
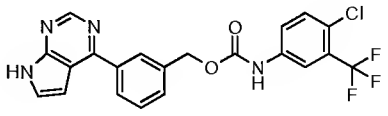
Entry	Name	Structure
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
133	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
134	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	

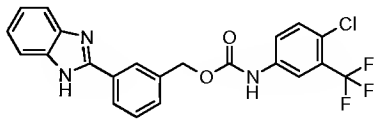
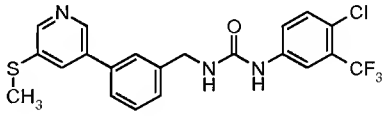
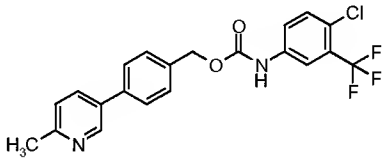
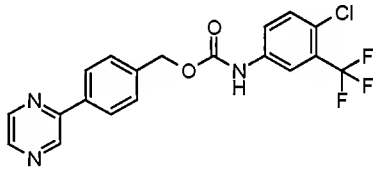
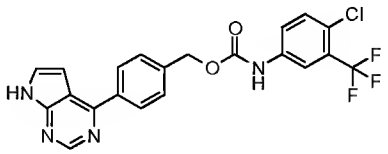
Entry	Name	Structure
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-indol-2-yl)phenyl]oxy}acetamide	

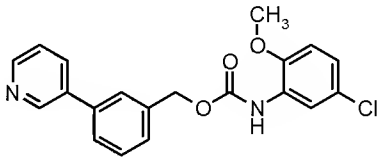
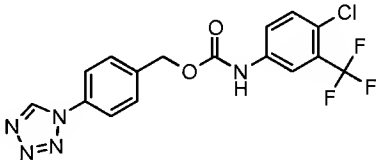
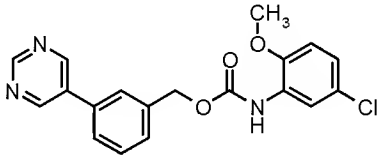
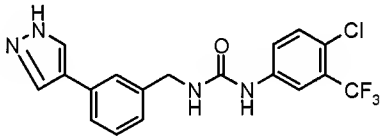
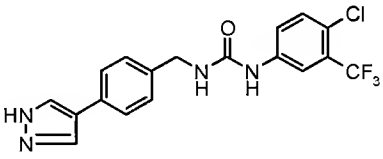
Entry	Name	Structure
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea	
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea	

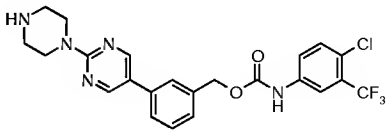
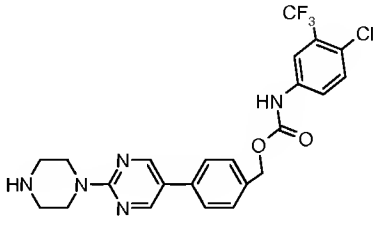
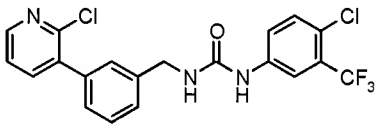
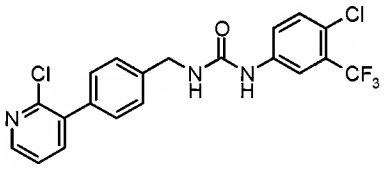
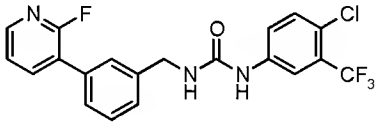
Entry	Name	Structure
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(5-hydroxy-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
160	N-{[3-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
161	N-{[4-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

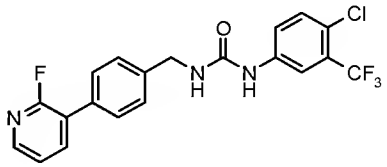
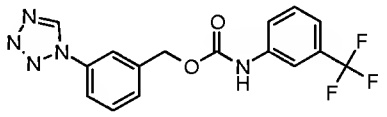
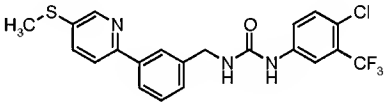
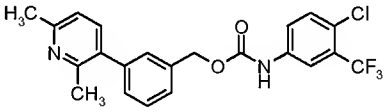
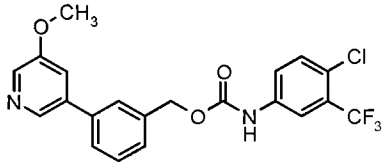
Entry	Name	Structure
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-methylpyridin-3-yl)phenyl]methyl}urea	

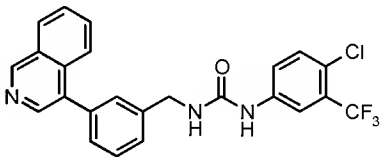
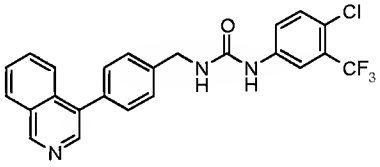
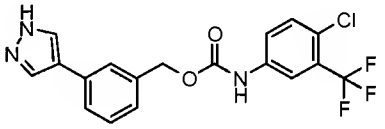
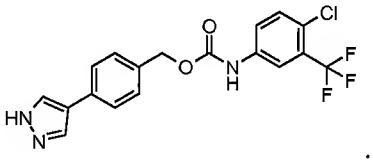
Entry	Name	Structure
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-methylpyridin-3-yl)phenyl]methyl}urea	
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
179	N-{[3-(6-acetylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-cyanopyridin-3-yl)phenyl]methyl}urea	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-3-yl]phenyl}methyl)urea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea	

Entry	Name	Structure
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N-{[3-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	

Entry	Name	Structure
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-2-yl]phenyl}methyl)urea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
226	{3-[5-(methyloxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-ylphenyl)methyl]urea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-ylphenyl)methyl]urea	
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (Previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.

29. (Cancelled)

30. (Withdrawn) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to claim 1.

31. (Withdrawn) The method according to claim 30, wherein the kinase is c-Kit.

32. (Withdrawn) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.

33. (Withdrawn) A method of treating rheumatoid arthritis, graft-host diseases, multiple sclerosis, psoriasis; artheroscrosis, myocardioinfarction, ischemia, stroke, restenosis; interbowel diseases, osteoarthritus, macular degeneration, or diabetic retinopathy, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 .

34. (Withdrawn) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.

35. (Withdrawn) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 to a cell or a plurality of cells.